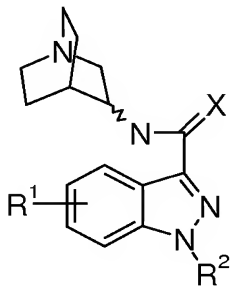
This listing of claims will replace all prior versions, and listings, of claims in the application:

Listing of Claims:

1. (Previously Presented): A compound of Formula I:



(I)

wherein

X is O or S;

R^1 is H, F, Cl, Br, I, OH, CN, nitro, NH_2, alkyl having 1 to 4 carbon atoms, fluorinated alkyl having 1 to 4 carbon atoms, cycloalkyl having 3 to 7 carbon atoms, cycloalkylalkyl having 4 to 7 carbon atoms, alkoxy having 1 to 4 carbon atoms, cycloalkoxy having 3 to 7 carbon atoms, cycloalkylalkoxy having 4 to 7 carbon atoms, alkylthio having 1 to 4 carbon atoms, fluorinated alkoxy having 1 to 4 carbon atoms, hydroxyalkyl having 1 to 4 carbon atoms, hydroxyalkoxy having 2 to 4 carbon atoms, monoalkylamino having 1 to 4 carbon atoms, dialkylamino wherein each alkyl group independently has 1 to 4 carbon atoms, Ar or Het;

R^2 is H, alkyl having 1 to 4 carbon atoms, cycloalkyl having 3 to 7 carbon atoms, or

cycloalkylalkyl having 4 to 7 carbon atoms;

Ar is an aryl group containing 6 to 10 carbon atoms which is unsubstituted or substituted one or more times by alkyl having 1 to 8 carbon atoms, alkoxy having 1 to 8 carbon atoms, halogen, dialkylamino wherein the alkyl portions each have 1 to 8 carbon atoms, amino, cyano, hydroxyl, nitro, halogenated alkyl having 1 to 8 carbon atoms, halogenated alkoxy having 1 to 8 carbon atoms, hydroxyalkyl having 1 to 8 carbon atoms, hydroxyalkoxy having 2 to 8 carbon atoms, alkenyloxy having 3 to 8 carbon atoms, alkylthio having 1 to 8 carbon atoms, alkylsulphinyl having 1 to 8 carbon atoms, alkylsulphonyl having 1 to 8 carbon atoms, monoalkylamino having 1 to 8 carbon atoms, cycloalkylamino wherein the cycloalkyl group has 3 to 7 carbon atoms and is optionally substituted, aryloxy wherein the aryl portion contains 6 to 10 carbon atoms and is optionally substituted, arylthio wherein the aryl portion contains 6 to 10 carbon atoms and is optionally substituted, cycloalkyloxy wherein the cycloalkyl group has 3 to 7 carbon atoms and is optionally substituted, sulfo, sulfonylamino, acylamido, acyloxy or combinations thereof; and

Het is a heterocyclic group, which is fully saturated, partially saturated or fully unsaturated, having 5 to 10 ring atoms in which at least 1 ring atom is a N, O or S atom, which is unsubstituted or substituted one or more times by halogen, aryl having 6 to 10 carbon atoms and is optionally substituted, alkyl having 1 to 8 carbon atoms, alkoxy having 1 to 8 carbon atoms, cyano, trifluoromethyl, nitro, oxo, amino, monoalkylamino having 1 to 8 carbon atoms, dialkylamino wherein each alkyl group has 1 to 8 carbon atoms, or combinations thereof; or

a pharmaceutically acceptable salt thereof,

wherein if the compound exhibits chirality it can be in the form of a mixture of enantiomers such as a racemate or a mixture of diastereomers, or can be in the form of a single

enantiomer or a single diastereomer.

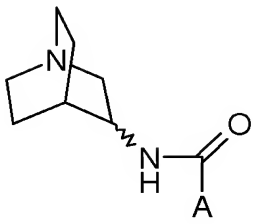
2. (Cancelled):

3. (Cancelled):

4. (Cancelled):

5. (Cancelled):

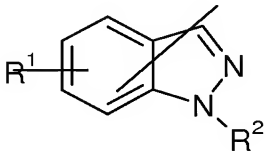
6. (Previously Presented): A compound according to Formula I':



(I')

wherein

A is an indazolyl according to subformula (a),

(a) 

R^1 is H, F, Cl, Br, I, OH, CN, nitro, NH_2, alkyl having 1 to 4 carbon atoms, fluorinated alkyl having 1 to 4 carbon atoms, cycloalkyl having 3 to 7 carbon atoms, cycloalkylalkyl having 4 to 7 carbon atoms, alkoxy having 1 to 4 carbon atoms, cycloalkoxy having 3 to 7 carbon atoms, alkylthio having 1 to 4 carbon

atoms, fluorinated alkoxy having 1 to 4 carbon atoms, hydroxyalkyl having 1 to 4 carbon atoms, hydroxyalkoxy having 2 to 4 carbon atoms, monoalkylamino having 1 to 4 carbon atoms, dialkylamino wherein each alkyl group independently has 1 to 4 carbon atoms, Ar or Het;

R^2 is H, alkyl having 1 to 4 carbon atoms, cycloalkyl having 3 to 7 carbon atoms, or cycloalkylalkyl having 4 to 7 carbon atoms;

Ar is an aryl group containing 6 to 10 carbon atoms which is unsubstituted or substituted one or more times by alkyl having 1 to 8 carbon atoms, alkoxy having 1 to 8 carbon atoms, halogen, dialkylamino wherein the alkyl portions each have 1 to 8 carbon atoms, amino, cyano, hydroxyl, nitro, halogenated alkyl having 1 to 8 carbon atoms, halogenated alkoxy having 1 to 8 carbon atoms, hydroxyalkyl having 1 to 8 carbon atoms, hydroxyalkoxy having 2 to 8 carbon atoms, alkenyloxy having 3 to 8 carbon atoms, alkylthio having 1 to 8 carbon atoms, alkylsulphinyl having 1 to 8 carbon atoms, alkylsulphonyl having 1 to 8 carbon atoms, monoalkylamino having 1 to 8 carbon atoms, cycloalkylamino wherein the cycloalkyl group has 3 to 7 carbon atoms and is optionally substituted, aryloxy wherein the aryl portion contains 6 to 10 carbon atoms and is optionally substituted, arylthio wherein the aryl portion contains 6 to 10 carbon atoms and is optionally substituted, cycloalkyloxy wherein the cycloalkyl group has 3 to 7 carbon ~~C~~ atoms and is optionally substituted, sulfo, sulfonylamino, acylamido, acyloxy or combinations thereof; and

Het is a heterocyclic group, which is fully saturated, partially saturated or fully unsaturated, having 5 to 10 ring atoms in which at least 1 ring atom is a N, O or S atom, which is unsubstituted or substituted one or more times by halogen, aryl having 6 to 10 carbon atoms and is optionally substituted, alkyl having 1 to 8 carbon atoms, alkoxy having 1 to 8 carbon atoms, cyano, trifluoromethyl, nitro,

oxo, amino, monoalkylamino having 1 to 8 carbon atoms, dialkylamino wherein each alkyl group has 1 to 8 carbon atoms, or combinations thereof; or

a pharmaceutically acceptable salt thereof,

wherein if the compound exhibits chirality it can be in the form of a mixture of enantiomers such as a racemate or a mixture of diastereomers, or can be in the form of a single enantiomer or a single diastereomer, and

wherein the indazolyl group of subformula (a) is attached to the remainder of the compound via its 3 position.

7. (Cancelled):

8. (Cancelled):

9. (Cancelled):

10. (Cancelled):

11. (Previously Presented): A compound according to claim 1, wherein R^1 is H, F, Cl, Br, 2-thiophenyl, 3-thiophenyl, 3-furyl, or phenyl.

12. (Cancelled):

13. (Cancelled):

14 (Previously Presented): A compound according to claim 1, wherein R^1 is H, F, Cl, Br, methyl, methoxy, or amino.

15. (Previously Presented): A compound according to claim 1, wherein R^2 is H or methyl.

16. (Cancelled):

17. (Cancelled):

18. (Cancelled):

19. (Cancelled):

20. (Cancelled):

21. (Previously Presented): A compound according to claim 1, wherein said compound is ~~selected fro~~m:

N-(1-Azabicyclo[2.2.2]oct-3-yl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof,

N-((3R)-1-Azabicyclo[2.2.2]oct-3-yl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof,

N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof,

1-Methyl-1H-Indazole-3-carboxamide, N-1-aza-bicyclo[2,2,2]oct-3-yl or a pharmaceutically acceptable salt thereof,

(R) 1-Methyl-1H-Indazole-3-carboxamide, N-1-aza-bicyclo[2,2,2]oct-3-yl or a pharmaceutically acceptable salt thereof,

(S) 1-Methyl-1H-Indazole-3-carboxamide, N-1-aza-bicyclo[2,2,2]oct-3-yl or a pharmaceutically acceptable salt thereof,

N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(bromo)-1H-indazole-3-carboxamide or a

pharmaceutically acceptable salt thereof,

N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(cyclopropyl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof,

N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(furan-3-yl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof,

N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(phenyl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof,

N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(thiophen-2-yl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof,

N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(thiophen-3-yl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof,

N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(bromo)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof,

N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(furan-3-yl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof,

N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(phenyl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof,

N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(thiophen-2-yl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof,

N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(thiophen-3-yl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof,

N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-6-(bromo)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof,

N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-6-(furan-3-yl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof,

N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-6-(phenyl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof,

N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-6-(thiophen-2-yl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof,

N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-6-(thiophen-3-yl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof,
N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-6-(bromo)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof,
N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-6-(furan-3-yl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof,
N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-6-(phenyl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof,
N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-6-(thiophen-2-yl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof,
N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-6-(thiophen-3-yl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof,
1-Aza-bicyclo[2,2,2]oct-3-ylcarboxamide, N-1H-indazol-3-yl or a pharmaceutically acceptable salt thereof,
(S) 1-Aza-bicyclo[2,2,2]oct-3-ylcarboxamide, N-1H-indazol-3-yl or a pharmaceutically acceptable salt thereof, or
(R) 1-Aza-bicyclo[2,2,2]oct-3-ylcarboxamide, N-1H-indazol-3-yl or a pharmaceutically acceptable salt thereof~~,~~
~~and pharmaceutically acceptable salts thereof~~.

22. (Previously Presented): A pharmaceutical composition comprising a compound according to claim 1 and a pharmaceutically acceptable carrier.

23. (Cancelled):

24. (Cancelled):

25. (Cancelled):

26. (Cancelled):

27. (Cancelled):

28. (Cancelled):

29. (Cancelled):

30. (Cancelled):

31. (Cancelled):

32. (Cancelled):

33. (Cancelled):

34. (Cancelled):

35. (Cancelled):

36. (Cancelled):

37. (Cancelled):

38. (Previously Presented): A compound according to claim 21, wherein said compound is in the form of a hydrochloride or hydroformate salt.

39. (Currently Amended): A compound according to claim 38, wherein said compound is ~~selected from~~:

N-(-1-Azabicyclo[2.2.2]oct-3-yl)-1H-indazole-3-carboxamide hydrochloride,
N-((3R)-1-Azabicyclo[2.2.2]oct-3-yl)-1H-indazole-3-carboxamide hydrochloride,
N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-1H-indazole-3-carboxamide hydrochloride,
N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(cyclopropyl)-1H-indazole-3-carboxamide hydroformate,
N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(furan-3-yl)-1H-indazole-3-carboxamide hydroformate,
N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(phenyl)-1H-indazole-3-carboxamide hydroformate,
N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(thiophen-2-yl)-1H-indazole-3-carboxamide hydroformate,
N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(thiophen-3-yl)-1H-indazole-3-carboxamide hydroformate,
N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(furan-3-yl)-1H-indazole-3-carboxamide hydroformate,
N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(phenyl)-1H-indazole-3-carboxamide hydroformate,
N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(thiophen-2-yl)-1H-indazole-3-carboxamide hydroformate,
N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(thiophen-3-yl)-1H-indazole-3-carboxamide hydroformate,
N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-6-(furan-3-yl)-1H-indazole-3-carboxamide hydroformate,
N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-6-(phenyl)-1H-indazole-3-carboxamide hydroformate,
N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-6-(thiophen-2-yl)-1H-indazole-3-carboxamide hydroformate,
N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-6-(thiophen-3-yl)-1H-indazole-3-carboxamide hydroformate,
N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-6-(furan-3-yl)-1H-indazole-3-carboxamide hydroformate,
N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-6-(phenyl)-1H-indazole-3-carboxamide hydroformate,
N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-6-(thiophen-2-yl)-1H-indazole-3-carboxamide hydroformate, or ~~and~~
N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-6-(thiophen-3-yl)-1H-indazole-3-carboxamide hydroformate.
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40. (Cancelled):

41. (Cancelled):

42. (Cancelled):

43. (Cancelled):

44. (Cancelled):

45. (Cancelled):

46. (Cancelled):

47. (Cancelled):

48. (Previously Presented): A compound according to claim 21, wherein said compound is N-(1-Azabicyclo[2.2.2]oct-3-yl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof.

49. (Previously Presented): A compound according to claim 48, wherein said compound is N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof.

50. (Previously Presented): A compound according to claim 48, wherein said compound is N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof.

51. (Previously Presented): A compound according to claim 21, wherein said

compound is N-(1-Azabicyclo[2.2.2]oct-3-yl)-5-(bromo)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof.

52. (Previously Presented): A compound according to claim 51, wherein said compound is N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(bromo)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof.

53. (Previously Presented): A compound according to claim 51, wherein said compound is N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(bromo)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof.

54. (Previously Presented): A compound according to claim 21, wherein said compound is N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(cyclopropyl)-1H-indazole-3-carboxamide carboxamide or a pharmaceutically acceptable salt thereof.

55. (Previously Presented): A compound according to claim 21, wherein said compound is N-(1-Azabicyclo[2.2.2]oct-3-yl)-5-(furan-3-yl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof.

56. (Previously Presented): A compound according to claim 55, wherein said compound is N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(furan-3-yl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof.

57. (Previously Presented): A compound according to claim 55, wherein said compound is N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(furan-3-yl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof.

58. (Previously Presented): A compound according to claim 21, wherein said compound is N-(1-Azabicyclo[2.2.2]oct-3-yl)-5-(phenyl)-1H-indazole-3-carboxamide or a

pharmaceutically acceptable salt thereof.

59. (Previously Presented): A compound according to claim 58, wherein said compound is N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(phenyl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof.

60. (Previously Presented): A compound according to claim 58, wherein said compound is N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(phenyl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof.

61. (Previously Presented): A compound according to claim 21, wherein said compound is N-(1-Azabicyclo[2.2.2]oct-3-yl)-5-(thiophen-2-yl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof.

62. (Previously Presented): A compound according to claim 61, wherein said compound is N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(thiophen-2-yl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof.

63. (Previously Presented): A compound according to claim 61, wherein said compound is N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(thiophen-2-yl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof.

64. (Previously Presented): A compound according to claim 21, wherein said compound is N-(1-Azabicyclo[2.2.2]oct-3-yl)-5-(thiophen-3-yl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof.

65. (Previously Presented): A compound according to claim 64, wherein said compound is N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(thiophen-3-yl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof.

66. (Previously Presented): A compound according to claim 64, wherein said compound is N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(thiophen-3-yl)-1H-indazole-3-carboxamide or a pharmaceutically acceptable salt thereof.

67. (Previously Presented): A compound according to claim 39, wherein said compound is N-(-1-Azabicyclo[2.2.2]oct-3-yl)-1H-indazole-3-carboxamide hydrochloride.

68. (Previously Presented): A compound according to claim 39, wherein said compound is N-((3R)-1-Azabicyclo[2.2.2]oct-3-yl)-1H-indazole-3-carboxamide hydrochloride.

69. (Previously Presented): A compound according to claim 39, wherein said compound is N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-1H-indazole-3-carboxamide hydrochloride.

70. (Previously Presented): A compound according to claim 39, wherein said compound is N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(cyclopropyl)-1H-indazole-3-carboxamide hydroformate.

71. (Previously Presented): A compound according to claim 39, wherein said compound is N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(furan-3-yl)-1H-indazole-3-carboxamide hydroformate.

72. (Previously Presented): A compound according to claim 39, wherein said compound is N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(phenyl)-1H-indazole-3-carboxamide hydroformate.

73. (Previously Presented): A compound according to claim 39, wherein said compound is N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(thiophen-2-yl)-1H-indazole-3-carboxamide hydroformate.

74. (Previously Presented): A compound according to claim 39, wherein said compound is N-((3*R*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(thiophen-3-yl)-1H-indazole-3-carboxamide hydroformate.

75. (Previously Presented): A compound according to claim 39, wherein said compound is N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(furan-3-yl)-1H-indazole-3-carboxamide hydroformate.

76. (Previously Presented): A compound according to claim 39, wherein said compound is N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(phenyl)-1H-indazole-3-carboxamide hydroformate.

77. (Previously Presented): A compound according to claim 39, wherein said compound is N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(thiophen-2-yl)-1H-indazole-3-carboxamide hydroformate.

78. (Previously Presented): A compound according to claim 39, wherein said compound is N-((3*S*)-1-Azabicyclo[2.2.2]oct-3-yl)-5-(thiophen-3-yl)-1H-indazole-3-carboxamide hydroformate.

79. (Previously Presented): A compound according to claim 11, wherein R^2 is H or methyl.

80. (Currently Amended): A compound according to claim 1, wherein Ar is substituted or unsubstituted phenyl or naphthyl, ~~naphthyl, or biphenyl,~~ and Het is substituted or unsubstituted tetrahydrofuranyl, tetrahydrothienyl, pyrrolidinyl, piperidinyl, piperazinyl, morpholinyl, isoxazolinyl, furyl, thienyl, pyrrolyl, pyrazolyl, imidazolyl, pyridyl, pyrimidinyl, indolyl, quinolinyl, isoquinolinyl, or naphthyridinyl.